L Number	Hits	Search Text	DB	Time stamp
1	0	("fluoren\$").PN.	USPAT;	2004/04/21 05:42
			US-PGPUB;	, , ,
			EPO; JPO;	
			DERWENT	
2	28987	fluoren\$	USPAT;	2004/04/21 05:42
1			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
3	2024	RAR	USPAT;	2004/04/21 05:42
			US-PGPUB;	
		·	EPO; JPO;	
	5257	63	DERWENT	
4	5257	fluorenone	USPAT;	2004/04/21 05:42
			US-PGPUB;	
	İ		EPO; JPO;	
5	0	RAR and fluorenone	DERWENT	2004/04/21 05 42
~		land tradictions	USPAT; US-PGPUB;	2004/04/21 05:42
· ·			EPO; JPO;	j
			DERWENT	
6	0	9807716.pn	USPAT;	2004/04/21 05:42
	1	F	US-PGPUB;	2004/04/21 05:42
			EPO; JPO;	j
		`	DERWENT	
7	586	562/466.ccls.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	*
	· ·		DERWENT	
8	189	549/26.ccls.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
9	108	514/437.ccls.	USPAT;	2004/04/21 05:42
	Į		US-PGPUB;	
			EPO; JPO;	ļ
10	524	514/569.ccls.	DERWENT	2004/04/05 05 10
10	524	514/569.CCIS.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO; DERWENT	
11	1331	562/466.ccls. or 549/26.ccls. or	USPAT;	2004/04/21 05:42
		514/437.ccls. or 514/569.ccls.	US-PGPUB;	2004/04/21 05.42
		,	EPO; JPO;	
·			DERWENT	
12	6541	retinoid	USPAT;	2004/04/21 05:42
			US-PGPUB;	, ,
		,	EPO; JPO;	
			DERWENT	
13	26,450	phenanth\$	USPAT;	2004/04/21 05:42
		•	US-PGPUB;	
			EPO; JPO;	
114	0441	rotinoia	DERWENT	0004/04/05 == -
14	8441	retinoic	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
15	113	retinoic and (562/466.ccls. or 549/26.ccls.	DERWENT USPAT;	2004/04/21 05:42
	113	or 514/437.ccls. or 514/569.ccls.)	US-PGPUB;	2004/04/21 05:42
	İ	,	EPO; JPO;	
		•	DERWENT	
16	853	560/56.ccls.	USPAT;	2004/04/21 05:42
		·	US-PGPUB;	, ,
			EPO; JPO;	
			DERWENT	
17	331	514/544.ccls.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	

18	2	5945561.pn.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
19	2	6319948.pn.	DERWENT	2004/24/23 25 10
19		6319946.pii.	USPAT;	2004/04/21 05:42
	1		US-PGPUB;	
			EPO; JPO;	
20	2	5760004 pp	DERWENT	0004/04/05 05 15
20	2	5760084.pn.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
21	2	FC240F7	DERWENT	
21	2	5624957.pn.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
20	4.5	63	DERWENT	
22	45	fluoren\$ and RAR	USPAT;	2004/04/21 05:42
			US-PGPUB;	1
1			EPO; JPO;	
	_		DERWENT	
23	5	9807716.pn.	USPAT;	2004/04/21 05:42
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
24	99	arotinoid	USPAT;	2004/04/21 05:43
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
25	1	arotinoid and 562/466.ccls.	USPAT;	2004/04/21 05:43
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
26	2	5618839.pn.	USPAT;	2004/04/21 05:43
		~	US-PGPUB;	,
			EPO; JPO;	
			DERWENT	
27	17	5618839.URPN.	USPAT	2004/04/21 05:43
28	3	arotinoid and (562/466.ccls. or 549/26.ccls.	USPAT;	2004/04/21 05:43
		or 514/437.ccls. or 514/569.ccls.)	US-PGPUB;	=====================================
			EPO; JPO;	1
			DERWENT	
29	139	retinoid and (562/466.ccls. or 549/26.ccls.	USPAT;	2004/04/21 05:43
		or 514/437.ccls. or 514/569.ccls.)	US-PGPUB;	-001/04/21 03.43
[-, · ·	EPO; JPO;	
			DERWENT	
30	4	(retinoid and (562/466.ccls. or 549/26.ccls.	USPAT;	2004/04/21 05:43
		or 514/437.ccls. or 514/569.ccls.)) and	US-PGPUB;	2004/04/21 03:43
		phenanth\$	EPO; JPO;	
[DERWENT	
31	3	phenanth\$ and (retinoic and (562/466.ccls.	USPAT;	2004/04/21 05:43
		or 549/26.ccls. or 514/437.ccls. or	US-PGPUB;	2004/04/21 05:43
		514/569.ccls.))	EPO; JPO;	
		J11/ JUJ. CC18. / /		
32	2	5075487.pn.	DERWENT	2004/04/21 25 42
22	- 4	2012±01.pm.	USPAT;	2004/04/21 05:43
			US-PGPUB;]
			EPO; JPO;	
33	853	560/56.ccls.	DERWENT	2004/04/21 26 25
33	853	500/50.CCIS.	USPAT;	2004/04/21 06:35
			US-PGPUB;	
	l		EPO; JPO;	
24	333	E14/E44 gglg	DERWENT	0004/0:/55
34	331	514/544.ccls.	USPAT;	2004/04/21 06:35
	ļ		US-PGPUB;	
			EPO; JPO;	
			DERWENT	

	Туре	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error	Definition
1	IS&R	L1	0	("fluoren\$").PN.	EPO; JPO; DERWE NT	2004/04/21 05:42			
2	BRS	L2	28987	fluoren\$		2004/04/21 05:42			
3	BRS	L3	2024	RAR	EPO; JPO; DERWE NT	2004/04/21 05:42			
4	BRS	L4	5257	fluorenone		2004/04/21 05:42			
5	BRS	L5	0	RAR and fluorenone	EPO; JPO; DERWE NT	2004/04/21 05:42			
6	BRS	L6	0	9807716.pn	EPO; JPO; DERWE NT	2004/04/21 05:42	·		
7	BRS	Ь7	586	562/466.ccls.		2004/04/21 05:42			

	Eri
1	0
2	0
3	0
4	0
5	0
6	0
7	0

	Туре	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error	Definition
8	BRS	L8	189	549/26.ccls.	EPO; JPO; DERWE NT	2004/04/21 05:42			
9	BRS	L 9	108	514/437.ccls.	EPO; JPO; DERWE NT	2004/04/21 05:42			
10	BRS	L10	524	514/569.ccls.	EPO; JPO; DERWE NT	2004/04/21 05:42			
11	BRS	L11	1331	562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.		2004/04/21 05:42			
12	BRS	L12	6541	retinoid	EPO; JPO; DERWE NT	2004/04/21 05:42			
13	BRS	L13	26450	phenanth\$	EPO; JPO; DERWE NT	2004/04/21 05:42			
14	BRS	L14	8441	retinoic		2004/04/21 05:42			

	Eri
8	0
9	0
10	0
11	0
12	0
13	0
14	0

	Туре	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error	Definition
15	BRS	L15	113	retinoic and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	EPO; JPO; DERWE NT	2004/04/21 05:42			
16	BRS	L16	853	560/56.ccls.	EPO; JPO; DERWE NT	2004/04/21 05:42			
17	BRS	L17	331	514/544.ccls.	EPO; JPO; DERWE NT	2004/04/21 05:42			
18	BRS	L18	2	5945561.pn.		2004/04/21 05:42			
19	BRS	L19	2	6319948.pn.		2004/04/21 05:42			
20	BRS	L20	2	5760084.pn.	EPO; JPO; DERWE NT	2004/04/21 05:42			-
21	BRS	L21	2	5624957.pn.		2004/04/21 05:42			

	Err
15	0
16	0
17	0
18	0
19	0
20	0
21	o

	Туре	L #	Hits	Search Text	DBs	Time	Stamp	Comments	Error	Defin	ition
22	BRS	L22	45	fluoren\$ and RAR	USPAT; US-PG PUB; EPO; JPO; DERWE	2004/ 05:42					
23	BRS	L23	5	9807716.pn.	EPO; JPO; DERWE NT	2004/ 05:42					
24	BRS	L24	99	arotinoid	EPO; JPO; DERWE NT	2004/ 05:43					
25	BRS	L25	1	arotinoid and 562/466.ccls.	USPAT; US-PG PUB; EPO; JPO; DERWE NT	2004/ 05:43				•	
26	BRS	L26	2	5618839.pn.	USPAT; US-PG PUB; EPO; JPO; DERWE NT	2004/ 05:43		١		*	
27	BRS	L27	17	5618839.URPN.	USPAT	2004/ 05:43	04/21				
28	BRS	L28	3 .	arotinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PG PUB; EPO; JPO; DERWE	2004/ 05:43					
29	BRS	L29	139	retinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PG PUB; EPO; JPO; DERWE	2004/ 05:43	/04/21 }				

22 0 23 0 24 0 25 0 26 0 27 0 28 0			Err
24 0 25 0 26 0 27 0 28 0	A CONTRACTOR OF THE PERSON OF	22	0
25 0 26 0 27 0 28 0		23	0
26 0 27 0 28 0		24	0
27 0		25	0
28 0		26	0
		27	0
29 0		28	0
		29	0

	Туре	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error	Definition
30	BRS	L30	4	(retinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)) and phenanth\$	EPO; JPO; DERWE NT	2004/04/21 05:43			
31	BRS	L31	3	phenanth\$ and (retinoic and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.))		2004/04/21 05:43			
32	BRS	L32	2	5075487.pn.	USPAT; US-PG PUB; EPO; JPO; DERWE NT	2004/04/21 05:43			,
33	BRS	L33	853	560/56.ccls.		2004/04/21 06:35			
34	BRS	L34	331	514/544.ccls.	USPAT; US-PG PUB; EPO; JPO; DERWE NT	2004/04/21 06:35			

	Err
30	0
31	0
32	0
33	0
34	0

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PASSWORD:

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=> FIL STNGUIDE

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

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FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Apr 16, 2004 (20040416/UP).

=> FIL HOME COST IN U.S. DOLLARS

SINCE FILE ENTRY TOTAL SESSION

FULL ESTIMATED COST

0.06

0.27

FILE 'HOME' ENTERED AT 05:54:10 ON 21 APR 2004

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

ENTRY

SESSION 0.48

0.21

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 19 APR 2004 HIGHEST RN 676225-08-4 DICTIONARY FILE UPDATES: 19 APR 2004 HIGHEST RN 676225-08-4

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=>

Uploading C:\Examination Auxillary files\10075845\10075845 generic six allowance.str

$$G_1$$
 G_1
 G_1
 G_2
 G_2
 G_2
 G_3
 G_1
 G_1
 G_1
 G_1
 G_2
 G_3
 G_4
 G_5
 G_6
 G_7
 chain nodes :
19 20 21 24 26 28 32
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 23 34
chain bonds :
10-26 13-24 16-19 19-20 19-21 20-32 24-26
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-34 7-8 7-9 8-12 8-23 9-10 10-11 11-12
13-18 13-14 14-15 15-16 16-17 17-18 23-34
exact/norm bonds :
5-7 6-34 8-23 10-26 13-24 16-19 19-20 19-21 20-32 23-34 24-26
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-9 8-12 9-10 10-11 11-12 13-18 13-14
14-15
15-16 16-17 17-18

G1:C,O,S,N

G2:C,O,S

G3:H,[*1]

Hydrogen count :
1:>= minimum 1 2:>= minimum 1 3:>= minimum 1 4:>= minimum 1
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 23:CLASS 24:CLASS 26:CLASS 28:CLASS 32:CLASS 34:CLASS Element Count :

Node 28: Limited C,C1-7

L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1

STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> search 11 sss sam
SAMPLE SEARCH INITIATED 05:54:49 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 3439 TO ITERATE

29.1% PROCESSED 1000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS:

65264 TO 72296

PROJECTED ANSWERS:

0 TO 0

L2 0 SEA SSS SAM L1

=> search 11 sss full FULL SEARCH INITIATED 05:55:01 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 69812 TO ITERATE

100.0% PROCESSED 69812 ITERATIONS SEARCH TIME: 00.00.03

0 ANSWERS

0 SEA SSS FUL L1

=>

L3

Uploading C:\Examination Auxillary files\10075845\10075845 generic fiveallowance.str

 $Ak^{\star 1}$ $28^{\star 1}$

chain nodes :
19 20 21 24 26 28 32
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 23
chain bonds :
10-26 13-24 16-19 19-20 19-21 20-32 24-26
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-23 7-8 7-9 8-12 8-23 9-10 10-11 11-12
13-18 13-14 14-15 15-16 16-17 17-18
exact/norm bonds :
5-7 6-23 8-23 10-26 13-24 16-19 19-20 19-21 20-32 24-26
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-9 8-12 9-10 10-11 11-12 13-18 13-14
14-15
15-16 16-17 17-18

G1:C,O,S,N

G2:C,O,S

G3:H, [*1]

L4 STRUCTURE UPLOADED

=> d 14 L4 HAS NO ANSWERS L4 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> search 14 sss sam
SAMPLE SEARCH INITIATED 05:57:01 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 3288 TO ITERATE

SAMPLE SCREEN SEARCH COMPLETED - 3288 TO ITERAT

30.4% PROCESSED 1000 ITERATIONS

0 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS:

62322 TO 69198

PROJECTED ANSWERS:

0 TO 0

L5 0 SEA SSS SAM L4

=> search 14 sss full FULL SEARCH INITIATED 05:57:09 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 66736 TO ITERATE

100.0% PROCESSED 66736 ITERATIONS

3 ANSWERS

SEARCH TIME: 00.00.03

L6 3 SEA SSS FUL L4

=> d scan

L6 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

MF C25 H21 N O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):3

L6 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Benzoic acid, 4-[(2-dibenzofuranylsulfonyl)amino]- (9CI)

MF C19 H13 N O5 S

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L6 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Benzoic acid, 4-[(2-dibenzofuranylsulfonyl)amino]-, methyl ester (9CI)
MF C20 H15 N O5 S

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 312.94 313.42

FULL ESTIMATED COST

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FILE COVERS 1907 - 21 Apr 2004 VOL 140 ISS 17 FILE LAST UPDATED: 20 Apr 2004 (20040420/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 16

L7

1 L6

=> d 17 ti fbib abs

L7 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN

TI Synthesis and biological activity of some new dibenzofuran- and 7-nitrodibenzofuran-2-sulfonyl amino acid derivatives

AN 1985:505304 CAPLUS

DN 103:105304

TI Synthesis and biological activity of some new dibenzofuran- and 7-nitrodibenzofuran-2-sulfonyl amino acid derivatives

AU El-Naggar, A. M.; Abd El-Salam, A. M.; Ahmed, F. S. M.; Ibrahim, T. M.

CS Fac. Sci., Al-Azhar Univ., Nasr, Egypt

SO Acta Pharmaceutica Jugoslavica (1985), 35(1), 15-22 CODEN: APJUA8; ISSN: 0001-6667

DT Journal

LA English

GΙ

$$\begin{array}{c|c} SO_2-X-OH & SO_2C1 \\ \hline I & R & O & III \\ \hline \\ R & O & III \\ \hline \end{array}$$

AB Title amino acid derivs. I (X = β -Ala, Val, Leu, p-NHC6H4CO, Phe, etc.; R = H or NO2) were prepared by sulfonylating the corresponding amino acid with sulfonyl chlorides II (R = H or NO2). I were esterified with MeOH via SOC12 to give the corresponding Me esters. Also, I were coupled with amino acid Me ester hydrochlorides by DCC in THF containing Et3N to give the corresponding dipeptides, e.g. III (X-X1 = DL-Val-DL-Val, Pro-Phe, R = H; X-X1 = Pro-DL-Ser, Leu-Tyr, R = NO2). Nineteen synthesized compds., e.g. I (X = Leu, R = H; X = β -Ala, R = NO2) and III (X-X1 = Tyr-Phe, R = NO2), were active against various microorganisms, e.g. Bacillus subtilis or B. cereus.

=> logoff hold COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 3.86 317.28 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

CA SUBSCRIBER PRICE

ENTRY -0.69

SESSION -0.69

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 06:00:04 ON 21 APR 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 06:06:12 ON 21 APR 2004 FILE 'CAPLUS' ENTERED AT 06:06:12 ON 21 APR 2004 COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS FULL ESTIMATED COST	SINCE FILE ENTRY 3.86	TOTAL SESSION 317.28
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) CA SUBSCRIBER PRICE	SINCE FILE ENTRY -0.69	TOTAL SESSION -0.69
=> file reg COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY 4.29	SESSION 317.71
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.69	-0.69

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STRUCTURE FILE UPDATES: 19 APR 2004 HIGHEST RN 676225-08-4 DICTIONARY FILE UPDATES: 19 APR 2004 HIGHEST RN 676225-08-4

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=> eBenzoic acid, 4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl ester/cn
MISSING OPERATOR '4-[[(9-OXO-9H-F'
```

=> e Benzoic ester/cn	acid,	4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl
E1	1	BENZOIC ACID, 4,6-DIMETHOXY-3-(5-METHOXY-2-(METHOXYCARBONYL) -3-PENTYLPHENOXY)-2-PENTYL-, METHYL ESTER/CN
E2	1	BENZOIC ACID, 4,6-DIMETHOXY-3-METHYL-2-(1-OXOBUTOXY)-, METHY L ESTER/CN
E3	0>	BENZOIC ACID, 4-(9-OXO-9H-FLUOREN-3-YL)CARBONYL AMINO-, 2 -METHYLPROPYL ESTER/CN
E4	1	BENZOIC ACID, 4-((((((((2,3-DIHYDRO-2,2-DIMETHYL-7-BENZOFURA NYL)OXY)CARBONYL)METHYLAMINO)CARBONYL)AMINO)SULFONYL)AMINO)- , ETHYL ESTER/CN
E5 .	1	BENZOIC ACID, 4-(((((((2,3,4-TRICHLOROPHENYL)THIO)ACETYL)OXY)ACETYL)AMINO)METHYL)-, METHYL ESTER/CN
E6	1	BENZOIC ACID, 4-(((((((3B)-17-OXOANDROST-5-EN-3-YL)AMIN O)CARBONYL)OXY)METHYL)AMINO)-, 2-(DIETHYLAMINO)ETHYL ESTER/CN
E7	1	BENZOIC ACID, 4-((((((3B)-17-OXOANDROST-5-EN-3-YL)OXY) METHYL)THIO)METHYL)AMINO)-, 2-(DIMETHYLAMINO)ETHYL ESTER/CN
E8	1	BENZOIC ACID, 4-(((((((3B,17B)-3-HYDROXYANDROST-5-EN-17-YL)AMINO)CARBONYL)OXY)METHYL)AMINO)-, 2-(DIETHYLAMINO)
Е9	1	ETHYL ESTER/CN BENZOIC ACID, 4-((((((4-BROMO-8-CHLORO-1-NAPHTHALENYL)THIO) ACETYL)OXY)ACETYL)AMINO)METHYL)-, METHYL ESTER/CN
E10	1	BENZOIC ACID, 4-((((((4-CHLOROPHENYL)AMINO)CARBONYL)HYDRAZO NO)PHENYLMETHYL)THIO)METHYL)-, 1,1-DIMETHYLETHYL ESTER/CN
E11	1	BENZOIC ACID, 4-(((((4-CHLOROPHENYL)SULFONYL)AMINO)PHENYLM ETHYLENE)AMINO)THIOXOMETHYL)AMINO)-/CN
E12	1	BENZOIC ACID, 4-(((((((4-METHOXYPHENYL)SULFONYL)(PHENYLMETHYL)AMINO)ACETYL)AMINO)OXY)DIPHENYLMETHYL)-/CN
=> e Benzoic ester/cn	acid,	4-(((9-oxo-9H-fluoren-3-yl)carbonyl)amino)-,2-methylpropyl
E1	1	BENZOIC ACID, 4-(((9-METHYL-4-OXO-2-(1-PYRROLIDINYL)-4H-PYRI DO(1,2-A)PYRIMIDIN-3-YL)METHYLENE)HYDRAZINO)-/CN
E2	1	BENZOIC ACID, 4-(((9-OXO-9H-FLUOREN-3-YL)CARBONYL)AMINO)-, 2 -METHYLPROPYL ESTER/CN
E3	0>	BENZOIC ACID, 4-(((9-OXO-9H-FLUOREN-3-YL)CARBONYL)AMINO)-,2-METHYLPROPYL ESTER/CN
E4	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-/CN
E5	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-, 1-(4-(2-AMINO-2-OXOETHOXY)-5-METHOXY-2-NITROPHENYL)ETHYL EST ER/CN
E6	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-, 1-METHYLETHYL ESTER/CN
E7	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-2-HYDROXY-/CN
E8	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-2-HYDROXY-, METHYL ESTER/CN
E9	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-2-NITRÒ-/CN
E10	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-3-IODO-/CN
E11	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-3-IODO-, METHYL ESTER/CN
E12	1	BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-3-METHOXY-/CN

=> e2

. Г8

1 "BENZOIC ACID, 4-(((9-OXO-9H-FLUOREN-3-YL)CARBONYL)AMINO)-, 2-METHYLPROPYL ESTER"/CN

=> d 18

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN

RN 505063-62-7 REGISTRY

CN Benzoic acid, 4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl ester (9CI) (CA INDEX NAME)

FS 3D CONCORD

MF C25 H21 N O4

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